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#### **MEMORANDUM**

DATE:

June 10, 2015

TO:

Eric Nuchims, Project Manager, E & E, Seattle, Washington

FROM:

Mark Woodke, START-4 Chemist, E & E, Seattle, Washington

SUBJ:

Organic Data Quality Assurance Review, John Day Vapor Response Site,

John Day, Oregon

REF:

TDD: 15-05-0005

PAN: 1004530.0004.111.02

The data quality assurance review of eight water samples collected from the John Day Vapor Response site in John Day, Oregon, has been completed. Semivolatile Organic Compound (SVOC) analysis (EPA Method 8270D SIM) was performed by TestAmerica, Inc., Tacoma, Washington. All sample analyses were evaluated following EPA's Stage 2B and/or 4 Data Validation Electronic and/or Manual Process (S2B/4VE/M).

The samples were numbered:

15053105

15053106

15053108

15053109

15053112

15053110

15053113

15053107

### **Data Qualifications:**

### 1. Sample Holding Times: Acceptable.

The samples were maintained and received within the QC limits of < 6°C. The samples were collected between May 29 and 31, 2015, were extracted on June 3, 2015, and were analyzed on June 6, 2015, therefore meeting holding time criteria of less than 7 days between collection and extraction and less than 40 days between extraction and analysis.

### 2. Tuning: Acceptable.

Tuning was performed at the beginning of each 12-hour analysis sequence. All results were within QC limits.

### 3. Initial Calibration: Acceptable.

All average Relative Response Factors (RRFs) were within the QC limits. All Relative Standard Deviations (RSDs) were within the QC limits.

### 4. Continuing Calibration: Acceptable.

All RRFs were within the QC limits. All % differences were within the QC limits.

# 5. Blanks: Acceptable.

A method blank was analyzed for each 20 sample batch per matrix. There were no detections in any method blank.

## 6. System Monitoring Compounds (SMCs): Acceptable.

All SMC recoveries were within QC limits.

# 7. Blank Spike (BS)/BS Duplicate (BSD) Analysis: Satisfactory.

All spike analyses were performed per SDG or per matrix per concentration level, whichever was more frequent. All recoveries were within the QC limits except acenaphthylene with low recoveries in the BS and BSD; associated positive results and sample quantitation limits were qualified as estimated with a low bias (JL or UJL).

## 8. Duplicate Analysis: Acceptable.

Blank spike duplicate analysis was performed per SDG or per matrix per concentration level, whichever was more frequent. All spike duplicate results were within QC limits.

## 9. Internal Standards: Acceptable.

All internal standards (IS) were within  $\pm$  30 seconds of the continuing calibration IS retention times. All area counts were within 50 % to 200 % of the continuing calibration area counts.

### 10. Precision and Bias Determination: Not Performed.

Samples necessary to determine precision and bias were not provided to the laboratory. All results were flagged "PND" (Precision Not Determined) and "RND" (Recovery Not Determined), although the flags do not appear on the data sheets.

### 11. Performance Evaluation Sample Analysis: Not Provided.

Performance evaluation samples were not provided to the laboratory.

#### 12. Overall Assessment of Data for Use

Pyrene (sample 15013105), phenanthrene (sample 15053106), and 2-methylnaphthalene (sample 15053107) were qualified as not detected (U) based on rinsate blank results.

The overall usefulness of the data is based on the criteria outlined in the Site-Specific Sampling Plan and/or Sampling and Quality Assurance Plan, the OSWER Guidance Document "Quality Assurance/Quality Control Guidance for Removal Activities, Sampling QA/QC Plan, and Data Validation Procedures" (EPA/540/G-90/004), the analytical method, and, when applicable, the Office of Emergency and Remedial Response Publication "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review". Based upon the information provided, the data are acceptable for use with the above stated data qualifications.

### Data Qualifiers and Definitions

- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- JH The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample with a high bias.
- JL The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample with a low bias.
- JK The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample with an unknown direction of bias.
- JQ The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample with an unknown direction of bias and falls between the MDL and the Minimum (or Practical) Quantitation Limit (MQL, PQL).
- N The analysis indicates the present of an analyte for which there is presumptive evidence to make a "tentative identification".
- NJ The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053105

Lab Sample ID:

580-50365-2

Client Matrix:

Water

Date Sampled: 05/29/2015 1216 Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method: 8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

580-191118

Lab File ID:

0606C011.D

Dilution:

Prep Batch:

Initial Weight/Volume:

1.0

Final Weight/Volume:

1055.5 mL 2.0 mL

Analysis Date: Prep Date:

06/06/2015 1943 06/03/2015 1647

Injection Volume:

1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Naphthalene	7.3		0.0068	0.019
2-Methylnaphthalene	6.4		0.0057	0.025
1-Methylnaphthalene	5.0		0.0057	0.019
Acenaphthylene	0.025 ブレ	*	0.0057	0.019
Acenaphthene	0.088		0.0057	0.019
Fluorene	0.14		0.0057	0.019
Phenanthrene	0.19		0.0057	0.019
Anthracene	<b>W</b> D		0.0057	0.019 📝
Fluoranthene	Mem		0.0057	0.019
Pyrene	0.022		0.0057	0.019
Benzo[a]anthracene	ND ND		0.0057	0.019()
Chrysene	, ND		0.0057	0.019
Benzo[b]fluoranthene	ήD		0.0057	0.019
Benzo[k]fluoranthene	ND		0.0057	0.019
Benzo[a]pyrene	ND		0.0057	0.019
Indeno[1,2,3-cd]pyrene	ND		0.0057	0.019
Dibenz(a,h)anthracene	иþ		0.0057	0.019
Benzo[g,h,i]perylene	ND		0.0057	0.019
	MV			
Surrogate	%Rec	Qualifier	Acceptar	ice Limits
Terphenyl-d14	94		64 - 150	~~~~

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053106

Lab Sample ID:

580-50365-3

Client Matrix:

Water

Date Sampled: 05/29/2015 1617

Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method: 8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

Prep Batch:

Lab File ID:

0606C012.D

Dilution:

1.0

580-191118

Initial Weight/Volume:

982.4 mL

Analysis Date:

Final Weight/Volume:

2.0 mL

Prep Date:

Surrogate

Terphenyl-d14

06/06/2015 2005 06/03/2015 1647

Injection Volume:

Qualifier

1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Naphthalene	0.30	***************************************	0.0073	0.020
2-Methylnaphthalene	0.18		0.0061	0.026
1-Methylnaphthalene	0.13		0.0061	0.020
Acenaphthylene	JAD	*	0.0061	0.020
Acenaphthene	NDING		0.0061	0.020()
Fluorene	0.0065	J <i>Q</i>	0.0061	0.020
Phenanthrene	0.0091	-JAM.	0.0061	0.020 <b>V</b> "
Anthracene	§ND	Ning	0.0061	0.020 🗸
Fluoranthene	ND		0.0061	0.020
Pyrene	ND		0.0061	0.020
Benzo[a]anthracene	ND		0.0061	0.020
Chrysene	ND		0.0061	0.020
Benzo[b]fluoranthene	Ν̈́D		0.0061	0.020
Benzo[k]fluoranthene	ND		0.0061	0.020
Benzo[a]pyrene	ΝD		0.0061	0.020
Indeno[1,2,3-cd]pyrene	ŊĎ.		0.0061	0.020
Dibenz(a,h)anthracene	ND		0.0061	0.020
Benzo[g,h,i]perylene	NP		0.0061	0.020
	. I Ba			

%Rec

98

MW.

Acceptance Limits

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053108

Lab Sample ID:

580-50365-4

Client Matrix:

Water

Date Sampled: 05/30/2015 1130

Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method: 8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

Prep Batch:

Lab File ID:

Dilution:

1.0

580-191118

0606C013.D

Initial Weight/Volume:

1054.1 mL

Analysis Date: Prep Date:

Terphenyl-d14

06/06/2015 2027 06/03/2015 1647 Final Weight/Volume: Injection Volume:

2.0 mL 1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL ,
Naphthalene	ND		0.0068	0.019
2-Methylnaphthalene	ND		0.0057	0.025
1-Methylnaphthalene	t <b>h</b> D	•	0.0057	0.019
Acenaphthylene	ΝD	*	0.0057	0.019
Acenaphthene	ΝD		0.0057	0.019
Fluorene	Ν̈́D		0.0057	0.019
Phenanthrene	ND		0.0057	0.019
Anthracene	ήνD		0.0057	0.019
Fluoranthene	ND		0.0057	0.019
Pyrene	ND		0.0057	0.019
Benzo[a]anthracene	ND		0.0057	0.019
Chrysene	ND		0.0057	0.019
Benzo[b]fluoranthene	ND		0.0057	0.019
Benzo[k]fluoranthene	ND		0.0057	0.019
Benzo[a]pyrene	ND		0.0057	0.019
Indeno[1,2,3-cd]pyrene	Мþ		0.0057	0.019
Dibenz(a,h)anthracene	NĎ		0.0057	0.019
Benzo[g,h,i]perylene	NI		0.0057	0.019
Surrogate	%Rec	Qualifier	Accepta	nce Limits

94

· WW Slate

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053109

Lab Sample ID:

580-50365-5

Client Matrix:

Water

Date Sampled: 05/30/2015 1428

Date Received: 06/02/2015 0945

#### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method:

8270D SIM 3520C

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

Prep Batch:

Result (ug/L)

580-191118

Lab File ID:

0606C014.D

Dilution:

1.0

Initial Weight/Volume:

MDL

0.0069

0.0057

0.0057

0.0057

0.0057

0.0057

1049.7 mL

Analysis Date: Prep Date:

06/06/2015 2049 06/03/2015 1647

Final Weight/Volume: Injection Volume:

Qualifier

2.0 mL 1 uL

RL

0.019

0.025

0.019

0.019

0.019

0.019

0.019

0.019

0.019

0.019

Analyte
Naphthalene
2-Methylnaphthalene
1-Methylnaphthalene
Acenaphthylene
Acenaphthene
Fluorene
Phenanthrene
Anthracene
Fluoranthene
Pyrene
Benzo[a]anthracene
Chrysene

ΝD ND ND ΝD

ΝD ND ΝD ΝD ΝD ND ΝD ΝD ND ND

0.0057 0.0057 0.0057 0.0057 0.0057 0.0057 0.0057 0.0057 0.0057

0.0057

0.0057

0.0057

0.019 0.019 0.019 0.019 0.019 0.019 0.019 0.019

Surrogate

Terphenyl-d14

Benzo[b]fluoranthene

Benzo[k]fluoranthene

Indeno[1,2,3-cd]pyrene

Dibenz(a,h)anthracene

Benzo[g,h,i]perylene

Benzo[a]pyrene

%Rec 98

Qualifier

Acceptance Limits

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053112

Lab Sample ID:

580-50365-6

Client Matrix:

Water

Date Sampled: 05/31/2015 1111

Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method:

8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

Lab File ID:

0606C015.D

Dilution:

1.0

Prep Batch:

580-191118

Initial Weight/Volume:

1048.6 mL 2.0 mL

Analysis Date:

06/06/2015 2110

Final Weight/Volume: Injection Volume:

Prep Date:

06/03/2015 1647

1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL 🚮
Naphthalene	ND		0.0069	0.019
2-Methylnaphthalene	ND		0.0057	0.025
1-Methylnaphthalene	ND		0.0057	0.019
Acenaphthylene	ND	*	0.0057	0.019 TL
Acenaphthene	ND		0.0057	0.019
Fluorene	ΝD		0.0057	0.019
Phenanthrene	ND	4	0.0057	0.019
Anthracene	ND		0.0057	0.019
Fluoranthene	ND		0.0057	0.019
Pyrene	ND		0.0057	0.019
Benzo[a]anthracene	ND		0.0057	0.019
Chrysene	ND		0.0057	0.019
Benzo[b]fluoranthene	Ν <mark></mark> D		0.0057	0.019
Benzo[k]fluoranthene	ND		0.0057	0.019
Benzo[a]pyrene	ND ND	•	0.0057	0.019
Indeno[1,2,3-cd]pyrene	N <b>þ</b> .		0.0057	0.019
Dibenz(a,h)anthracene	NID		0.0057	0.019
Benzo[g,h,i]perylene	ND		0.0057	0.019 <b>V</b>
Surrogate	%Rec	Qualifier	Acceptance Limits	
Terphenyl-d14	90	64 - 150		

MW HOB

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053110

Lab Sample ID:

580-50365-7

Client Matrix:

Water

Date Sampled: 05/30/2015 1748

Date Received: 06/02/2015 0945

# 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method:

8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

Prep Batch:

Lab File ID:

0606C016.D

Dilution:

1.0

580-191118

Initial Weight/Volume:

1055.8 mL

Analysis Date:

06/06/2015 2132

Final Weight/Volume: Injection Volume:

2.0 mL

06/03/2015 1647

1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL /
Naphthalene	ND	titi italan enatrini itari eksistitii ita iraanan (n. 1911).	0.0068	0.019 🗸
2-Methylnaphthalene	ND		0.0057	0.025
1-Methylnaphthalene	ND		0.0057	0.019
Acenaphthylene	ŊD	*	0.0057	0.019
Acenaphthene	ND		0.0057	0.019
Fluorene	ΝID		0.0057	0.019
Phenanthrene	ND		0.0057	0.019
Anthracene ·	Ν̈́D		0.0057	0.019
Fluoranthene	ND		0.0057	0.019
Pyrene	ND		0.0057	0.019
Benzo[a]anthracene	ND		0.0057	0.019
Chrysene	ND		0.0057	0.019
Benzo[b]fluoranthene	ND		0.0057	0.019
Benzo[k]fluoranthene	ND		0.0057	0.019
Benzo[a]pyrene	ΝĎ		0.0057	0.019
Indeno[1,2,3-cd]pyrene	ND		0.0057	0.019
Dibenz(a,h)anthracene	ND		0.0057	0.019
Benzo[g,h,i]perylene	ND		0.0057	0.019
Surrogate	%Rec	Qualifier	Acceptance Limits	
Terphenyl-d14	96	64 - 150		

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053113

Lab Sample ID:

580-50365-8

Client Matrix:

Water

Date Sampled: 05/31/2015 1433 Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method:

8270D SIM

Analysis Batch: 580-191434

Instrument ID:

TAC023

Prep Method:

3520C

Lab File ID:

0606C017.D

Dilution:

1.0

Prep Batch:

580-191118

Initial Weight/Volume:

1054.3 mL

Analysis Date:

06/06/2015 2154

Final Weight/Volume: Injection Volume:

2.0 mL 1 uL

Prep Date:

Terphenyl-d14

06/03/2015 1647

Analyte	Result (ug/L)	Qualifier	MDL	RL 11
Naphthalene	ND		0.0068	0.019 🗸
2-Methylnaphthalene	ND .		0.0057	0.025
1-Methylnaphthalene	ND .		0.0057	0.019
Acenaphthylene	ND ·	*	0.0057	0.019
Acenaphthene	ŅD		0.0057	0.019
Fluorene	ΝD		0.0057	0.019
Phenanthrene	Ν̈́D		0.0057	0.019
Anthracene	ΝD		0.0057	0.019
Fluoranthene	ŊD		0.0057	0.019
Pyrene	ND		0.0057	0.019
Benzo[a]anthracene	ND	*	0.0057	0.019
Chrysene	ŊD		0.0057	0.019
Benzo[b]fluoranthene	ND		0.0057	0.019
Benzo[k]fluoranthene	ŊD		0.0057	0.019
Benzo[a]pyrene	ND		0.0057	0.019
Indeno[1,2,3-cd]pyrene	Nþ .		0.0057	0.019
Dibenz(a,h)anthracene	ΝΦ		0.0057	0.019
Benzo[g,h,i]perylene	ND		0.0057	0.019
Surrogate	%Rec	Qualifier	Acceptar	nce Limits

91

CMW KIDIS

Client: Ecology and Environment, Inc.

Job Number: 580-50365-1

Client Sample ID:

15053107

Lab Sample ID:

580-50365-9

Client Matrix:

Water

Date Sampled: 05/31/2015 1100

Date Received: 06/02/2015 0945

### 8270D SIM Semivolatile Organic Compounds (GC/MS SIM)

Analysis Method:

8270D SIM 3520C

Analysis Batch:

580-191434

Instrument ID:

**TAC023** 

Prep Method:

Prep Batch:

Lab File ID:

0606C018.D

Dilution:

1.0

580-191118

Initial Weight/Volume:

979.7 mL

Analysis Date: Prep Date:

06/06/2015 2216 06/03/2015 1647

Final Weight/Volume: Injection Volume:

2.0 mL 1 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Naphthalene	0.014	ΙQ	0.0073	0.020
2-Methylnaphthalene	- <del>0.00</del> 91		0.0061	0.027 <b>U</b>
1-Methylnaphthalene	ND	. `	0.0061	0.020
Acenaphthylene	Ν̈́D	*	0.0061	0.020 TTC
Acenaphthene	ND		0.0061	0.020
Fluorene	ND		0.0061	0.020
Phenanthrene	ŊD		0.0061	0.020
Anthracene	ND		0.0061	0.020
Fluoranthene	Nb		0.0061	0.020
Pyrene	ΝΦ		0.0061	0.020
Benzo[a]anthracene	Иф		0.0061	0.020
Chrysene	ND		0.0061	0.020
Benzo[b]fluoranthene	ΝΦ		0.0061	0.020
Benzo[k]fluoranthene	ΝΦ		0.0061	0.020
Benzo[a]pyrene	ND		0.0061	0.020
Indeno[1,2,3-cd]pyrene	ND	•	0.0061	0.020
Dibenz(a,h)anthracene	ND		0.0061	0.020
Benzo[g,h,i]perylene	ND		0.0061	0.020
Surrogate	%Rec	Qualifier	er Acceptance Limits	
Terphenyl-d14	97	. 64 - 150		